#### (FILE 'HOME' ENTERED AT 15:26:42 ON 14 DEC 2005)

L31

```
FILE 'REGISTRY' ENTERED AT 15:26:50 ON 14 DEC 2005
L1
             26 S C39H38N2O2/MF
             98 S C33H40N2O2/MF
L2
             16 S L2 AND AMINO
L3
L4
              0 S L2 AND PHENOL
L5
              3 S L2 AND DIMETHYLPHENYL
L6
             6 S L2 AND (HYDROXY (3W) PHENYL)
L7
            16 S L2 AND AMINO
L8
            133 S C33H38N2O2/MF
L9
            34 S L8 AND AMINO
L10
            133 S C33H38N2O2/MF
     FILE 'CAPLUS' ENTERED AT 15:46:43 ON 14 DEC 2005
L11
              6 S (620951-13-5# OR 651306-73-9# OR 651306-75-1#)/RN
     FILE 'REGISTRY' ENTERED AT 15:53:54 ON 14 DEC 2005
L12
           1051 S METHYLENEBIS AND AMINO AND PHENOL
L13
             1 S 22445-96-1/RN
L14
             17 S C13H14N2O2/MF AND AMINO AND PHENOL
L15
             82 S METHYLENEBIS AND BENZYL AND AMINO
L16
             7 S METHYLENEBIS AND BENZYL AND AMINO AND PHENOL
L17
           3834 S AMINOPHENOL
L18
           443 S 3 (W) AMINOPHENOL
           645 S 4 (W) AMINOPHENOL
L19
L20
             61 S C20H21N2O2/MF
Ĺ21
             0 S L20 AND METHYLENBIS AND AMINO AND PHENOL
L22
            19 S L20 AND AMINO
L23
           1051 S METHYLENEBIS AND AMINO AND PHENOL
L24
            1 S 22445-96-1/RN
L25
           4090 S PHENOL AND METHYLETHYLIDENE AND AMINO
L26
             1 S 63969-46-0/RN
L27
              1 S 46999-95-5/RN
L28
           336 S PHENOL AND METHYLETHYLIDENE (W) BIS (3W) AMINO
L29
            61 S C20H21N2O2/MF
L30
             0 S L29 AND METHYLENEBIS AND AMINO AND PHENOL
```

0 S L29 AND METHYLENEBIS AND AMINO AND ((HYDROXY(3W)PHENYL###)) OR

#### (FILE 'HOME' ENTERED AT 17:04:50 ON 14 DEC 2005)

```
FILE 'REGISTRY' ENTERED AT 17:04:58 ON 14 DEC 2005
L1
             0 S TRIS (3W) HYDROXYPHENOL (3W) METHANE
             26 S TRIS (3W) HYDROXYPHENYL (3W) METHANE
L2
L3
              9 S 123-30-8/CRN AND 50-00-0/CRN AND 108-95-2/CRN
              1 S 95-55-6/CRN AND 50-00-0/CRN AND 108-95-2/CRN
L4
              1 S 62-53-3
L5
L6
            116 S 62-53-3/CRN AND 50-00-0/CRN AND 108-95-2/CRN
              2 S 62-53-3/CRN AND 50-00-0/CRN AND 90-15-3/CRN
L7
L8
              1 S 62-53-3/CRN AND 50-00-0/CRN AND 135-19-3/CRN
L9
             51 S 62-53-3/CRN AND 50-00-0/CRN AND 80-05-7/CRN
             0 S 62-53-3/CRN AND 50-00-0/CRN AND 603-44-1/CRN
L10
L11
             94 S (BISPHENOL (W) F) (3A) (DIGLYCIDYL (W) ETHER)
     FILE 'CAPLUS' ENTERED AT 17:14:21 ON 14 DEC 2005
          36157 S (1675-54-3# OR 25085-99-8# OR 25068-38-6# OR 65581-98-8# OR 2
L12
            426 S (24937-74-4# OR 38891-71-3# OR 38806-75-6#)/RN
L13
                S EP/PCT
     FILE 'REGISTRY' ENTERED AT 17:16:24 ON 14 DEC 2005
L14
          47901 S EP/PCT
     FILE 'CAPLUS' ENTERED AT 17:16:24 ON 14 DEC 2005
L15
         73288 S L14
L16
         176763 S (EPOX!###(W)RESIN#) OR DIEPOX!### OR POLYEPOX!### OR DIGLYCID
=> s 112 or 115 or 116
L17 186826 L12 OR L15 OR L16
=> s l17 and l13
L18
         162 L17 AND L13
```

```
65581-98-8 REGISTRY
RN
ED
     Entered STN: 16 Nov 1984
     Oxirane, 2,2'-[methylenebis(4,1-phenyleneoxymethylene)]bis-, homopolymer
CN
           (CA INDEX NAME)
     (9CI)
OTHER NAMES:
     A 5040
CN
     Araldite GY 282
CN
     Araldite GY 285
CN
CN
     Araldite LY 5082
CN
     Araldite PY 306
CN
     Beckopox SEP 106
CN
     Bisphenol F diglycidyl ether homopolymer
CN
     Epalloy 8220
CN
     EPC 830
     Epiclon 830
CN
     Epiclon 830CR
CN
CN
     Epiclon EXA 830CRP
CN
     Epiclon EXA 830LVP
     Epikote 806
CN
    Epikote 862
CN
     EXA 830CRP
CN
CN
     EXA 830LVP
     LY 5082
CN
     PY 306
CN
CN
     SBF 19
     SBF 22
CN
CN
     SBF 24
CN
     Vantico GY 282
CN
     YL 6475
     YL 983L
CN
     448927-09-1, 177646-13-8, 130501-68-7, 133516-56-0, 154214-00-3,
DR
     159777-68-1, 188797-94-6
MF
     (C19 H20 O4)x
CI
     PMS, COM
PCT
     Epoxy resin, Polyether
LC
                  AGRICOLA, CA, CAPLUS, PIRA, PROMT, TOXCENTER, USPAT2,
     STN Files:
       USPATFULL
     CM
          1
     CRN 2095-03-6
     CMF C19 H20 O4
```

377 REFERENCES IN FILE CA (1907 TO DATE)
73 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
377 REFERENCES IN FILE CAPLUS (1907 TO DATE)

```
2095-03-6 REGISTRY
RN
ED
     Entered STN: 16 Nov 1984
     Oxirane, 2,2'-[methylenebis(4,1-phenyleneoxymethylene)]bis- (9CI)
CN
                                                                          (CA
     INDEX NAME)
OTHER CA INDEX NAMES:
    Methane, bis[p-(2,3-epoxypropoxy)phenyl]- (7CI, 8CI)
CN
OTHER NAMES:
     4,4'-Methylenebisphenol diglycidyl ether
CN
     4,4'-Methylenediphenol diglycidyl ether
CN
CN
     Bis (4-glycidyloxyphenyl) methane
CN
     Bis (4-hydroxyphenyl) methane diglycidyl ether
CN
     Bis (p-hydroxyphenyl) methane diglycidyl ether
CN
    Bisphenol F diglycidyl ether
CN
     Diphenylolmethane diglycidyl ether
    GY 281
CN
CN
    p,p-BFDGE
CN
     para-para-BFDGE
FS
     3D CONCORD
MF
     C19 H20 O4
CI
     COM
LC
     STN Files: AGRICOLA, ANABSTR, BEILSTEIN*, BIOBUSINESS, BIOSIS, CA,
       CAOLD, CAPLUS, CASREACT, CBNB, CHEMCATS, CHEMLIST, CSCHEM, CSNB, IFICDB,
       IFIPAT, IFIUDB, PIRA, RTECS*, TOXCENTER, USPAT7ULL
         (*File contains numerically searchable property data)
     Other Sources:
                      EINECS**
         (**Enter CHEMLIST File for up-to-date regulatory information)
```

202 REFERENCES IN FILE CA (1907 TO DATE)

54 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA

202 REFERENCES IN FILE CAPLUS (1907 TO DATE)

3 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

RN 39817-09-9 REGISTRY

ED Entered STN: 16 Nov 1984

CN Oxirane, 2,2'-[methylenebis(phenyleneoxymethylene)]bis- (9CI) (CA INDEX NAME)

OTHER CA INDEX NAMES:

CN Methane, bis[(2,3-epoxypropoxy)phenyl] - (7CI)

OTHER NAMES:

CN Bisphenol F bis(oxiranylmethyl) ether

CN Bisphenol F diglycidyl ether

DR 87110-76-7

MF C19 H20 O4

CI IDS, COM

LC STN Files: AGRICOLA, BIOBUSINESS, BIOSIS, CA, CAOLD, CAPLUS, CHEMLIST, MSDS-OHS, PIRA, TOXCENTER, USPAT2, USPATFULL

Other Sources: EINECS\*\*, NDSL\*\*, TSCA\*\*

(\*\*Enter CHEMLIST File for up-to-date regulatory information)



- 133 REFERENCES IN FILE CA (1907 TO DATE)
- 62 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
- 133 REFERENCES IN FILE CAPLUS (1907 TO DATE)
  - 1 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

```
RN
     24937-74-4 REGISTRY
ED
     Entered STN: 16 Nov 1984
     Formaldehyde, polymer with benzenamine and phenol (9CI)
                                                                (CA INDEX NAME)
CN
OTHER CA INDEX NAMES:
     Aniline, polymer with formaldehyde and phenol (8CI)
CN
     Benzenamine, polymer with formaldehyde and phenol (9CI)
CN
CN
     Phenol condensation products, with aniline and formaldehyde (8CI)
     Phenol, polymer with benzenamine and formaldehyde (9CI)
CN
OTHER NAMES:
CN
     AF 10
     AF 10 (polymer)
CN
CN
     AFF 214
     AFFS 214
CN
     Aniline-formaldehyde-phenol copolymer
CN
     Aniline-formaldehyde-phenol polymer
CN
CN
     Aniline-formaldehyde-phenol resin
     Aniline-formaldehyde-phenolic copolymer
CN
CN
     Aniline-phenol-formaldehyde copolymer
CN
     Aniline-phenol-formaldehyde polymer
CN
     Aniline-phenol-formaldehyde resin
CN
     FAF
CN
     FAF (binder)
CN
     FAF (resin)
CN
     Formaldehyde-aniline-phenol resin
CN
     K 211-2
CN
     K 211-3
     K 211-34
CN
     K 214-2
CN
     K 214-22
CN
CN
     Phenol formaldehyde aniline polymer
CN
     Phenol-aniline-formaldehyde polymer
CN
     Phenol-aniline-formaldehyde resin
CN
     Phenol-formaldehyde-aniline copolymer
CN
     Phenol-formaldehyde-aniline resin
CN
     R-ANP
CN
     RAFF 214
CN
     Resin 211
CN
     Resin 214
CN
     Resol 214
CN
     RSM 143T
CN
     SF 340
CN
     SF 340A
     SF 341A
CN
CN
     SF 342
CN
     SF 342A
CN
     SVAF 2
CN
     Voloknit STV 1
DR
     9041-59-2, 9082-39-7, 9082-40-0, 149659-75-6
MF
     (C6 H7 N . C6 H6 O . C H2 O)x
CI
     PMS
PCT
     Amino resin, Phenolic resin
LC
                 BIOSIS, CA, CAPLUS, CHEMLIST, IFICDB, IFIPAT, IFIUDB,
     STN Files:
       NIOSHTIC, PIRA, PROMT, TOXCENTER, USPATFULL
     Other Sources:
                      DSL**, TSCA**
         (**Enter CHEMLIST File for up-to-date regulatory information)
     CM
          1
          108-95-2
     CRN
     CMF
          C6 H6 O
```

CM 2

CRN 62-53-3 CMF C6 H7 N

CM :

CRN 50-00-0 . CMF C H2 O

 $H_2C = 0$ 

\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

402 REFERENCES IN FILE CA (1907 TO DATE)

32 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA

402 REFERENCES IN FILE CAPLUS (1907 TO DATE)

```
RN
     38891-71-3 REGISTRY
```

ED

Entered STN: 16 Nov 1984
Formaldehyde, polymer with benzenamine and 1-naphthalenol (9CI) (CA INDEX NAME)

OTHER CA INDEX NAMES:

1-Naphthalenol, polymer with benzenamine and formaldehyde (9CI) CN

Benzenamine, polymer with formaldehyde and 1-naphthalenol (9CI) CN

OTHER NAMES:

Aniline-formaldehyde- $\alpha$ -naphthol polymer CN

MF (C10 H8 O . C6 H7 N . C H2 O) x

CI PMS

Amino resin, Phenolic resin PCT

LC STN Files: CA, CAPLUS, IFICDB, IFIPAT, IFIUDB, USPATFULL

CM

90-15-3 CRN CMF C10 H8 O

CM 2

CRN 62-53-3 CMF C6 H7 N

CM 3

CRN 50-00-0 CMF C H2 O

H2C=0

- 1 REFERENCES IN FILE CA (1907 TO DATE)
- 1 REFERENCES IN FILE CAPLUS (1907 TO DATE)

RN 38806-75-6 REGISTRY

ED

Entered STN: 16 Nov 1984
Formaldehyde, polymer with benzenamine and 4,4'-(1-CN methylethylidene)bis[phenol] (9CI) (CA INDEX NAME)

OTHER CA INDEX NAMES:

Benzenamine, polymer with formaldehyde and 4,4'-(1methylethylidene)bis[phenol] (9CI)

Phenol, 4,4'-(1-methylethylidene)bis-, polymer with benzenamine and CN formaldehyde (9CI)

OTHER NAMES:

Aniline-bisphenol A-formaldehyde copolymer CN

Aniline-bisphenol A-formaldehyde polymer CN

CN Aniline-formaldehyde-bisphenol A copolymer

(C15 H16 O2 . C6 H7 N . C H2 O)  $\times$ MF

CI PMS

PCT Amino resin, Phenolic resin

LC STN Files: CA, CAPLUS, IFICDB, IFIPAT, IFIUDB, USPATFULL

CM 1

CRN 80-05-7 CMF C15 H16 O2

CM 2

CRN 62-53-3 CMF C6 H7 N

CM 3

CRN 50-00-0 CMF C H2 O

 $H_2C = 0$ 

30 REFERENCES IN FILE CA (1907 TO DATE)

3 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA

30 REFERENCES IN FILE CAPLUS (1907 TO DATE)

RN 620951-13-5 REGISTRY

ED Entered STN: 26 Nov 2003

CN Phenol, 2,2'-methylenebis[6-[(4-amino-3,5-dimethylphenyl)methyl]-4-

methyl- (9CI) (CA INDEX NAME)

OTHER NAMES:

CN 2,2'-Methylenebis[6-(4-amino-3,5-dimethylbenzyl)-4-methylphenol]

FS 3D CONCORD

MF C33 H38 N2 O2

CI COM

SR CA

LC STN Files: CA, CAPLUS, USPATFULL

spec., p. 14, Diamine

- 6 REFERENCES IN FILE CA (1907 TO DATE)
- 3 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
- 6 REFERENCES IN FILE CAPLUS (1907 TO DATE)

RN 651306-73-9 REGISTRY

ED Entered STN: 18 Feb 2004

CN 2-Naphthalenol, 1-[(4-amino-3,5-dimethylphenyl)methyl]-3-[[3-[(4-amino-3,5-dimethylphenyl)methyl]-2-hydroxy-1-naphthalenyl]methyl]- (9CI) (CA INDEX

MF C39 H38 N2 O2

CI COM

SR CA

LC STN Files: CA, CAPLUS, USPATFULL

$$\begin{array}{c} \text{Me} \\ \text{Me} \\ \text{CH}_2 \\ \text{HO} \\ \text{CH}_2 \\ \text{HO} \\ \text{CH}_2 \\ \text{Me} \\ \text{CH}_2 \\ \text{Me} \\ \text{Me} \\ \text{CH}_2 \\$$

spec., p. 14, DISMINE 2

- 2 REFERENCES IN FILE CA (1907 TO DATE)
- 1 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
- 2 REFERENCES IN FILE CAPLUS (1907 TO DATE)

RN 651306-75-1 REGISTRY

ED Entered STN: 18 Feb 2004

CN Phenol, 2,6-bis[(4-amino-3,5-dimethylphenyl)methyl]-4-[1-(4-hydroxyphenyl)-1-methylethyl]- (9CI) (CA INDEX NAME)

FS 3D CONCORD

MF C33 H38 N2 O2

CI COM

SR CA

LC STN Files: CA, CAPLUS, USPATFULL

spec., p. 14, Diamine 3

- 2 REFERENCES IN FILE CA (1907 TO DATE)
- 1 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
- 2 REFERENCES IN FILE CAPLUS (1907 TO DATE)

RN 63969-46-0 REGISTRY

ED Entered STN: 16 Nov 1984

CN Phenol, 2,2'-methylenebis[4-amino- (6CI, 9CI) (CA INDEX NAME)

OTHER NAMES:

CN Bis(5-amino-2-hydroxyphenyl)methane

FS 3D CONCORD

MF C13 H14 N2 O2

CI COM

LC STN Files: CA, CAOLD, CAPLUS, IFICDB, IFIPAT, IFIUDB, TOXCENTER, USPATFULL

- 62 REFERENCES IN FILE CA (1907 TO DATE)
- 1 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
- 62 REFERENCES IN FILE CAPLUS (1907 TO DATE)
- 1 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

RN 16523-28-7 REGISTRY

ED Entered STN: 16 Nov 1984

CN Phenol, 4,4'-methylenebis[2-amino- (7CI, 8CI, 9CI) (CA INDEX NAME)

OTHER NAMES:

CN 3,3'-Diamino-4,4'-dihydroxydiphenylmethane

CN 4,4'-Methylenebis (2-aminophenol)

CN Bis (3-amino-4-hydroxyphenyl) methane

FS 3D CONCORD

MF C13 H14 N2 O2

CI COM

LC STN Files: BEILSTEIN\*, CA, CAOLD, CAPLUS, CHEMCATS (\*File contains numerically searchable property data)

- 16 REFERENCES IN FILE CA (1907 TO DATE)
- 2 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
- 16 REFERENCES IN FILE CAPLUS (1907 TO DATE)
- 1 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

RN 22428-30-4 REGISTRY

ED Entered STN: 16 Nov 1984

CN Phenol, 3,3'-methylenebis[6-amino- (7CI, 8CI, 9CI) (CA INDEX NAME)

OTHER NAMES:

CN 3,3'-Dihydroxy-4,4'-diaminodiphenylmethane

CN 4,4'-Diamino-3,3'-dihydroxydiphenylmethane

FS 3D CONCORD

MF C13 H14 N2 O2

CI COM

LC STN Files: BEILSTEIN\*, CA, CAOLD, CAPLUS, CHEMCATS, DETHERM\*, IFICDB, IFIUDB, TOXCENTER, USPATFULL

(\*File contains numerically searchable property data)

$$H_2N$$
 OH OH

- 16 REFERENCES IN FILE CA (1907 TO DATE)
- 2 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
- 16 REFERENCES IN FILE CAPLUS (1907 TO DATE)
- 1 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

RN 6423-19-4 REGISTRY

ED Entered STN: 16 Nov 1984

CN Phenol, 4,4'-methylenebis[3-amino- (8CI, 9CI) (CA INDEX NAME)

FS 3D CONCORD

MF C13 H14 N2 O2

CI COM

LC STN Files: CA, CAOLD, CAPLUS

# \*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

5 REFERENCES IN FILE CA (1907 TO DATE)

5 REFERENCES IN FILE CAPLUS (1907 TO DATE)

1 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

```
1220-78-6 REGISTRY
RN
ED
     Entered STN: 16 Nov 1984
     Phenol, 4,4'-(1-methylethylidene)bis[2-amino-(9CI) (CA INDEX
     NAME)
OTHER CA INDEX NAMES:
     Phenol, 4,4'-isopropylidenebis[2-amino- (7CI, 8CI)
CN
OTHER NAMES:
     2,2-Bis(3-amino-4-hydroxyphenyl)propane
CN
     4,4'-Isopropylidenebis(2-aminophenol)
CN
     DAM 1 (diol)
CN
CN
     NSC 10847
    3D CONCORD
FS
DR
     27905-27-7
MF
     C15 H18 N2 O2
CI
     COM
LC
     STN Files:
                  BEILSTEIN*, CA, CAOLD, CAPLUS, CASREACT, CHEMCATS, IFICDB,
       IFIPAT, IFIUDB, TOXCENTER, USPAT2, USPATFULL
         (*File contains numerically searchable property data)
```

- 36 REFERENCES IN FILE CA (1907 TO DATE)
- 5 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
- 37 REFERENCES IN FILE CAPLUS (1907 TO DATE)
- 4 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

RN 38361-73-8 REGISTRY

ED Entered STN: 16 Nov 1984

CN Phenol, 4,4'-(1-methylethylidene)bis[2-amino-6-methyl- (9CI)

(CA INDEX NAME)

OTHER CA INDEX NAMES:

CN o-Cresol, 4,4'-isopropylidenebis[6-amino- (7CI)

OTHER NAMES:

CN 2,2-Bis(3-amino-4-hydroxy-5-methylphenyl)propane

CN 2,2-Bis(3-methyl-4-hydroxy-5-aminophenyl)propane

FS 3D CONCORD

MF C17 H22 N2 O2

CI COM

LC STN Files: CA, CAOLD, CAPLUS

- 5 REFERENCES IN FILE CA (1907 TO DATE)
- 2 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
- 5 REFERENCES IN FILE CAPLUS (1907 TO DATE)
- 2 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

RN 22445-96-1 REGISTRY

ED Entered STN: 16 Nov 1984

CN Phenol, 3,3'-(1-methylethylidene)bis[6-amino-(9CI) (CA INDEX NAME)

OTHER CA INDEX NAMES:

CN Phenol, 3,3'-isopropylidenebis[6-amino- (8CI)

FS 3D CONCORD

MF C15 H18 N2 O2

CI COM

$$\begin{array}{c|c} & \text{Me} \\ \hline \\ \text{H}_2\text{N} & \text{OH} \\ \end{array}$$

\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

63969-A6-D

```
RN
     123-30-8 REGISTRY
ED
     Entered STN: 16 Nov 1984
     Phenol, 4-amino- (9CI) (CA INDEX NAME)
CN
OTHER CA INDEX NAMES:
     Phenol, p-amino- (8CI)
CN
OTHER NAMES:
     1-Amino-4-hydroxybenzene
CN
     4-Amino-1-hydroxybenzene
CN
CN
     4-Aminophenol
     4-Hydroxy-1-aminobenzene
CN
CN
     4-Hydroxyaniline
CN
     4-Hydroxybenzenamine
CN
     4-Hydroxyphenylamine
CN
     Activol
CN
     Azol
     BASF Ursol P Base
CN
     Benzofur P
CN
CN
     C.I. 76550
     C.I. Oxidation Base 6
CN
CN
     Certinal
CN
     Citol
CN
     Durafur Brown RB
     Fouramine P
CN
CN
     Fourrine 84
     Fourrine P Base
CN
CN
     Furro P base
     Nako Brown R
CN
CN
     NSC 1545
CN
     p-Aminophenol
CN
     p-Hydroxyaniline
CN
     p-Hydroxyphenylamine
CN
     Paranol
CN
     Pelagol Grey P Base
CN
     Pelagol P Base
CN
     Renal AC
CN
     Rodinal
     Tertral P Base
CN
CN
     Unal
CN
     Ursol P
CN
     Ursol P Base
CN
     Zoba Brown P Base
FS
     3D CONCORD
DR
     52985-09-8
MF
     C6 H7 N O
CI
LC
                  ADISNEWS, AGRICOLA, ANABSTR, AQUIRE, BEILSTEIN*, BIOBUSINESS,
       BIOSIS, BIOTECHNO, CA, CANCERLIT, CAOLD, CAPLUS, CASREACT, CBNB, CEN,
       CHEMCATS, CHEMINFORMRX, CHEMLIST, CIN, CSCHEM, CSNB, DDFU, DETHERM*,
       DRUGU, EMBASE, GMELIN*, HODOC*, HSDB*, IFICDB, IFIPAT, IFIUDB,
       IMSCOSEARCH, IPA, MEDLINE, MRCK*, MSDS-OHS, NIOSHTIC, PIRA, PROMT, PS,
       RTECS*, SCISEARCH, SPECINFO, SYNTHLINE, TOXCENTER, ULIDAT, USPAT2,
       USPATFULL
         (*File contains numerically searchable property data)
                      DSL**, EINECS**, TSCA**
         (**Enter CHEMLIST File for up-to-date regulatory information)
```

7543 REFERENCES IN FILE CA (1907 TO DATE)

513 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA

7557 REFERENCES IN FILE CAPLUS (1907 TO DATE)

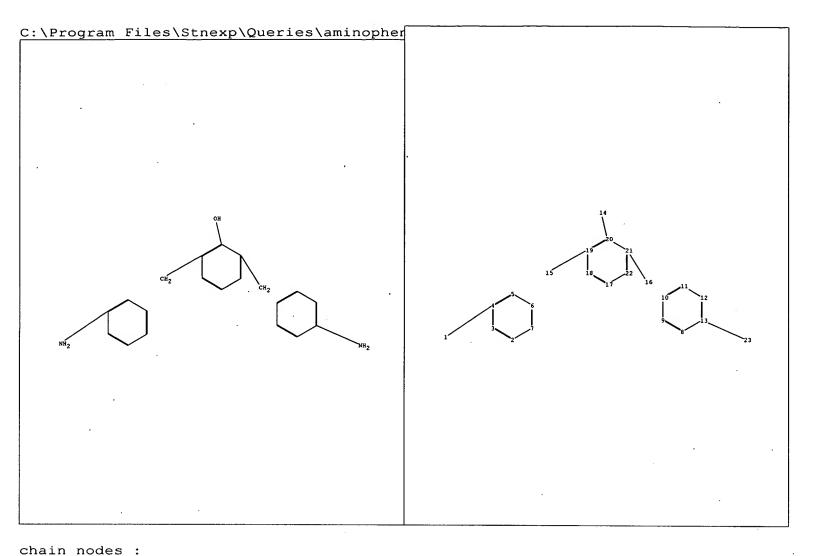
8 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

```
95-55-6 REGISTRY
RN
ED
     Entered STN: 16 Nov 1984
     Phenol, 2-amino- (9CI) (CA INDEX NAME)
CN
OTHER CA INDEX NAMES:
     Phenol, o-amino- (8CI)
CN
OTHER NAMES:
     (2-Hydroxyphenyl)amine
CN
CN
     1-Amino-2-hydroxybenzene
CN
     1-Hydroxy-2-aminobenzene
     2-Amino-1-hydroxybenzene
CN
CN
     2-Aminophenol
     2-Hydroxyaniline
CN
CN
     2-Hydroxybenzenamine
CN
     BASF Ursol 3GA
     Benzofur GG
CN
     C.I. 76520
CN
CN
     C.I. Oxidation Base 17
     Fouramine OP
CN
     Nako Yellow 3GA
CN
     NSC 1534
CN
     NSC 226261
CN
     o-Aminohydroxybenzene
CN
CN
     o-Aminophenol
     o-Hydroxyaniline
CN
     o-Hydroxyphenylamine
CN
     Paradone Olive Green B
CN
CN
     Pelagol 3GA
CN
     Pelagol Grey GG
CN
     Rodol 2G
CN
     Zoba 3GA
FS
     3D CONCORD
     C6 H7 N O
MF
CI
     COM
LC
     STN Files:
                  AGRICOLA, ANABSTR, AQUIRE, BEILSTEIN*, BIOBUSINESS, BIOSIS,
       BIOTECHNO, CA, CANCERLIT, CAOLD, CAPLUS, CASREACT, CHEMCATS,
       CHEMINFORMRX, CHEMLIST, CIN, CSCHEM, CSNB, DDFU, DETHERM*, DRUGU,
       EMBASE, GMELIN*, HODOC*, HSDB*, IFICDB, IFIPAT, IFIUDB, IPA, MEDLINE,
       MRCK*, MSDS-OHS, NIOSHTIC, PIRA, PROMT, PS, RTECS*, SPECINFO, SYNTHLINE,
       TOXCENTER, ULIDAT, USPAT2, USPATFULL
         (*File contains numerically searchable property data)
                      DSL**, EINECS**, TSCA**
         (**Enter CHEMLIST File for up-to-date regulatory information)
```

6064 REFERENCES IN FILE CA (1907 TO DATE)
344 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
6080 REFERENCES IN FILE CAPLUS (1907 TO DATE)
5 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

```
RN
     591-27-5 REGISTRY
ED
     Entered STN: 16 Nov 1984
     Phenol, 3-amino- (9CI) (CA INDEX NAME)
CN
OTHER CA INDEX NAMES:
     Phenol, m-amino- (8CI)
OTHER NAMES:
     (3-Hydroxyphenyl)amine
CN
     1-Amino-3-hydroxybenzene
CN
CN
     3-Aminophenol
CN
     3-Hydroxyaniline
     3-Hydroxybenzenamine
CN
     C.I. 76545
CN
     C.I. Oxidation Base 7
CN
     Fouramine EG
CN
CN
     Fourrine 65
     Fourrine EG
CN
     Furro EG
CN
     Futramine EG
CN
CN
     m-Aminophenol
CN
     m-Hydroxyaminobenzene
CN
     m-Hydroxyaniline
     m-Hydroxyaniline
CN
     m-Hydroxyphenylamine
CN
CN
     Nako TEG
     NSC 1546
CN
     Pelagol EG
CN
CN
     Renal EG
CN
     Rodol EG
CN
     Tertral EG
CN
     Ursol EG
CN
     Zoba EG
FS
     3D CONCORD
MF
     C6 H7 N O
CI
     COM
                  ADISNEWS, AGRICOLA, ANABSTR, AQUIRE, BEILSTEIN*, BIOBUSINESS,
LC
     STN Files:
       BIOSIS, BIOTECHNO, CA, CANCERLIT, CAOLD, CAPLUS, CASREACT, CBNB, CEN,
       CHEMCATS, CHEMINFORMRX, CHEMLIST, CIN, CSCHEM, CSNB, DDFU, DETHERM*,
       DRUGU, EMBASE, GMELIN*, HODOC*, HSDB*, IFICDB, IFIPAT, IFIUDB, IPA,
       MEDLINE, MRCK*, MSDS-OHS, NIOSHTIC, PIRA, PROMT, PS, RTECS*, SPECINFO,
       SYNTHLINE, TOXCENTER, ULIDAT, USPAT7, USPATFULL
         (*File contains numerically searchable property data)
                      DSL**, EINECS**, TSCA**
         (**Enter CHEMLIST File for up-to-date regulatory information)
```

4215 REFERENCES IN FILE CA (1907 TO DATE)
328 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
4226 REFERENCES IN FILE CAPLUS (1907 TO DATE)
4 REFERENCES IN FILE CAOLD (PRIOR TO 1967)



```
ring nodes :
    2 3 4 5 6 7
                          8 9 10 11 12 13 17 18
                                                             19
                                                                  20
                                                                       21
chain bonds :
    1-4 13-23 14-20 15-19 16-21
ring bonds:

2-3 2-7 3-4 4-5 5-6 6-7 8-9 8-13 9-10 10-11 11-12 12-13 17-18 17-22 18-19 19-20 20-21 21-22
exact/norm bonds :
     1-4 13-23 14-20
exact bonds :
    15-19 16-21
normalized bonds :
     2-3 2-7 3-4 4-5 5-6 6-7 8-9 8-13 9-10 10-11
                                                                      11-12 12-13 17-18
     17-22 18-19 19-20 20-21 21-22
Match level:
    1:CLASS 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:Atom 11:Atom 12:Atom 13:Atom 14:CLASS 15:CLASS 16:CLASS 17:Atom 18:Atom 19:Atom 20:Atom 21:Atom 22:Atom 23:CLASS
```

1 14 15 16 23